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New methods for accelerating the determination of basis trees and dual evaluators for distribution problems are compared with standard solution procedures in a computational study of a wide range of distribution problems of varying sizes and densities. Computer programs utilizing the new methods are tested for computational efficiency in an experiment involving four solution techniques, four start algorithms, and four change of basis criteria, thus affording an empirical determination not only of the merits of various procedures in isolation but also of their effectiveness in combination.

The study discloses that the most efficient solution procedure arises by coupling a primal transportation algorithm (embodying the accelerated updating and pricing methods) with a version of the Row Minimum start rule and a "modified first negative evaluator" rule. The resulting method was found to improve upon the efficiency of general purpose algorithms (taken from standard computer packages) by a factor of 50 or better, and also improved upon a streamlined version of the SHARE out-of-kilter code by a factor of 3. The method's median solution time for solving 175 x 175 distribution problems on a CDC 6600 computer was 11.4 seconds with a range of 9 to 13 seconds.

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A COMPARISON OF COMPUTATION TIMES FOR VARIOUS STARTING PROCEDURES, BASIS CHANGE CRITERIA, AND SOLUTION ALGORITHMS FOR DISTRIBUTION PROBLEMS

by

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May 1971

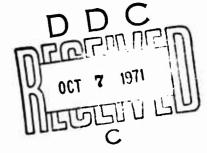
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## **ABSTRACT**

New methods for accelerating the determination of basis trees and dual evaluators for distribution problems are compared with standard solution procedures in a computational study of a wide range of distribution problems of varying sizes and densities. Computer programs utilizing the new methods are tested for computational efficiency in an experiment involving four solution techniques, four start algorithms, and four change of basis criteria, thus affording an empirical determination not only of the merits of various procedures in isolation but also of their effectiveness in combination.

The study discloses that the most efficient solution procedure arises by coupling a primal transportation algorithm (embodying the accelerated updating and pricing methods) with a version of the Row Minimum start rule and a "modified first negative evaluator" rule. The resulting method was found to improve upon the efficiency of general purpose algorithms (taken from standard computer packages) by a factor of 50 or better, and also improved upon a streamlined version of the SHARE out-of-kilter code by a factor of 3. The method's median solution time for solving 175 x 175 distribution problems on a CDC 6600 computer was 11.4 seconds with a range of 9 to 13 seconds.

# 1.0 Introduction and Scope of the Computational Analysis

The major purpose of this paper is to examine the following notions that have become part of the folklore surrounding the distribution (transportation) problem. 1

- (1) The most efficient algorithm for solving network problems is the out-of-kilter network algorithm developed by Ford and Fulkerson [11];
- (2) The computational efficiency provided by special purpose algorithms for distribution and network problems is not significantly greater than that provided by general purpose linear programming methods that use sophisticated procedures for exploiting sparse matrices;
- (3) The most efficient criterion (from the standpoint of total computational time) for determining a basis exchange in the context of the distribution problem is the most negative evaluator criterion [9, 20, 21];
- (4) The total solution time required to solve distribution problems of dimension m x m increases proportionately to  $m^2$  [10].

A secondary purpose of this study is to compare solution times between special purpose primal and dual simplex algorithms for solving distribution problems. Such a comparison is motivated by the conjecture [18, 19] that the special purpose dual algorithms are likely to be "faster" than their counterpart primal algorithms. The reason for this conjecture is two-fold: (a) To apply the most negative evaluator criterion to a dense m x m distribution problem in the dual approach requires the examination of only (2m-1) evaluators; whereas the primal approach requires the examination of m<sup>2</sup> - (2m-1) evaluators; (b) there exist procedures for finding basic feasible solutions for the dual of

the distribution problem while counterpart procedures (which insure that the starting basis will always be feasible) do not exist for the primal problem. Another purpose of our study, related to the foregoing, is to examine the effect on total computational time of different procedures for obtaining initial starting bases for both the primal and dual formulations of the distribution problem. The final purpose of our study is to examine the effect of the density on the solution time of an m x m distribution problem.

To guarantee a valid and impartial comparison of the procedures under analysis, the distribution problems used in the study varied between 9 percent and 100 percent density and varied in size from 10 x 10 to 200 x 200 (origins x destinations). A total of 140 different distribution problems were examined, all of which were randomly generated using a uniform probability function. The total supply of each m x m distribution problem was set equal to 1000m. The only other restrictions placed on the problems consisted of requiring the number of variables to be less than or equal to 10,000 and requiring the cost coefficients to lie between 1 and 100.

To accomplish these purposes, it has been necessary to obtain (and in some instances to develop) computer codes that are representative of the "state of the art" in solving network and linear programming problems. A description of the codes used in this study (e.g., out-of-kilter algorithm, general linear programming algorithm, and special purpose primal and dual simplex algorithms for solving distribution problems) is contained in a subsequent section.

The CDC 6600 at The University of Texas at Austin Computation Center was used to solve the distribution problems for the out-of-kilter and epecial purpose primal and dual simplex computer codes. Furthermore, the computer jobs were executed during periods when the machine load was approximately the same. The CDC 6600 at the CDC Data Center in Houston was used to solve the distribution problems for the general purpose simplex computer code, OPHELIE/LP.

The practical importance of determining the efficiency of alternative ways for solving distribution and network problems is affirmed not only by the fact that a sizeable fraction of the linear programming literature has been devoted to it, but also by the fact that an even larger share of the many concrete industrial and military applications of linear programming deal with distribution problems. Distribution problems often occur as subproblems in a larger problem (e.g., the traveling salesman problem or the warehouse location problem). Moreover, industrial applications of distribution problems often contain thousands of variables, and hence a streamlined algorithm is not only computationally worthwhile but a practical necessity. In addition, a number of linear programs that appear at first glance to be unrelated to the distribution problem can nevertheless be given a distribution problem formulation [7, 9, 11, 22, 23, 33], and it is also possible to approximate certain additional linear programming problems by such a formulation.

# 2.0 Description of Procedures Under Analysis

# 2.1 Solution Algorithms

The general simplex linear programming computer code employed in the study was Control Data's OPHELIE/LP code. OPHELIE/LP is a subsystem of the OPHELIE II Mathematical Programming System, and it fully exploits the characteristics of the CDC 6600 computer.

The optimization is performed by a primal algorithm based on the revised simplex method with the product form of the inverse. The algorithm performs the computation with a selection of multiple columns that is not fixed a priori but that can be altered dynamically to take into account statistics established in the course of iterations and the space available in central memory.

Two alternatives are permitted: one may successively introduce the best columns from among the remaining candidates, or optimize the linear sub-program formed by the basis and candidates. If cycling is detected during solution, an automatic perturbation procedure is triggered to remedy it. For further information concerning OPHELIE/LP, see [26, 27].

The other algorithms used in the study are designed to take direct advantage of the topological structure of the distribution problem.

The special purpose primal simplex approach is the "Row-Column Sum Method" [7]. To enhance the computational efficiency of the algorithm, the predecessor and augmented predecessor index methods [17, 18] are employed to determine a basis exchange path and to effect the appropriate updating of the problem data. (The basis for these procedures is the triple labeling

procedure proposed by Ellis Johnson [23].)

Lemke's Dual Method [25] provides the foundation for the special purpose dual algorithm<sup>3</sup> (the "Dual Row-Column Sum Method"). The computer code for this algorithm modifies the dual method to take advantage of the predecessor and augmented predecessor index methods and the poly-ω method [7] to produce what has been called the "double pricing" dual algorithm [15]. Both the primal and dual row-column methods are "in core" computer codes. Each currently has a limit of 10,000 admirsible cells (variables).

The out-of-kilter algorithm was tested both because of its reputation as a highly efficient solution method for network problems in general and also because of the sparseness of some of the distribution problems used in this study (which the out-of-kilter method is capable of exploiting to special advantage). The computer code for the out-of-kilter method employed in the analysis was written by R. J. Classen of the RAND Corporation and is distributed by SHARE [8, 30]. The out-of-kilter computer code is an "in core" algorithm and currently has a limit of 10,000 admissible arcs. The general form of the code was modified slightly by the authors to enhance its computational efficiency for application to distribution problems.

# 2.2 Start Algorithms

A variety of claims have been made concerning the influence of various "starting" algorithms, although documented evidence to support these claims has not been entirely satisfactory. To help remedy this situation, our study includes comparisons of several primal "start" methods that have

appeared in the literature. The primal start methods examined are the Northwest Corner Rule [9, 20], Vogel A pproximation Method (VAM) [7, 31], Row-Column Minimum [9], and Row-Minimum [10, 20]. These start algorithms do not exhaust all of those that have been proposed, but they are representative of the start procedures available for the primal formulation of the distribution problem.

The dual start method examined (which is of more recent vintage than the primal start methods) is described in [15, 19].

# 2.3 Basis Change Criteria

Four different criteria for specifying the variable to enter the basis (in the primal approach) were examined to ascertain their effect upon the total solution time: the commonly proposed "first negative evaluator" [10] and "most negative evaluator" [9] criteria, and variations of these called the "modified first negative evaluator" [10] and "modified most negative evaluator" criteria. The "first negative evaluator" scans the rows of the distribution tableau until it encounters the first element whose evaluator signals that this cell may "profitably" enter the basis. The "modified first negative evaluator" criterion scans the rows of the distribution tableau until it encounters the first row that contains a negative (i.e., profitable) evaluator, and then selects the most negative evaluator in this row.

The "modified most negative method" begins by selecting the most negative evaluator in the tableau on the initial iteration of the algorithm. At each subsequent iteration of the algorithm, a satisficing criterion is used that

selects the first variable encountered which has an evaluator less than some multiple of the evaluator selected in the preceding iteration. If no such variable exists, the method selects the variable with the most negative evaluator. This criterion is highly flexible, since it can be made to select the same variable as the "first negative evaluator" criterion by using a zero multiple of the preceding evaluator and to select the same variable as the "most negative evaluator" criterion by using a large multiple of the preceding evaluator. The multiple used in our analysis was 1/3 (developed by the authors on the basis of preliminary testing).

In the primal approach it is important to consider the trade-off between the amount of time required to implement various basis change criteria and the number of iterations that result by using these criteria, since a method that results in fewer iterations may still consume greater overall time due to an increased time per iteration.

In the dual approach such a trade-off is not particularly important for dense m x m distribution problems since, instead of considering up to m<sup>2</sup> - (2m-1) evaluators (as in the primal approach), the dual method has to contend with only 2m-1 evaluators. Consequently, only the most negative basis change criterion (which tends to reduce the total number of iterations) was used in the special purpose dual simplex code.

# 3.0 Numerical Results and Their Interpretation

In this section we present the results of the computational experiment concerning folklores, start algorithms, solution algorithms and basis

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change criteria. The various median computation times appear in Tables 1-5.

The computation times do not include the time for input and output.

# 3.1 Folklores

The results of the analysis indicate that the topological structure of the distribution model merits the use of special purpose linear programming algorithms to solve such problems. The data in Tables 1 and 2 indicate that the primal and dual row-column sum algorithms are, respectively, at least 20 times faster than the OPHELIE/LP general simplex linear programming computer code. In general the special purpose primal simplex method is 50 times faster. For instance, the solution for a 100 x 100 distribution is approximately 4.5 seconds using the special purpose code and 277 seconds using the OPHELIE/LP code. (Before computing the median times for the OPHELIE/LP code, a number of trial runs were made to determine the best procedure for selecting the next incoming cell. The trial results indicated that 30 cells should be considered at a time and, further, the most negative cell should be selected.)

Furthermore, the analysis indicates that the out-of-kilter algorithm is not the most efficient solution technique for solving all network problems. Even for extremely sparse distribution problems, as indicated in Tables 1 and 2, the special purpose primal algorithm, using the row-minimum start method and the modified first negative basis change criterion, was consistently 2 to 3 times faster than the out-of-kilter algorithm. Unexpectedly, as

Table 1

Solution Times (sec) for Out-of-Kilter, OPHELIE/LP and Dual Algorithms

Time per	. 002	. 005	600.	.010	. 020	.015	. 023	.030	.034	. 030
Total IT*	.017	. 202	. 639	. 845	2.947	2,929	6.873	9,633	14,956	14.387
Dual <sup>1</sup> No. Iterations	ω	37	70	83	144	201	293	324	446	476
Start	. 008	. 024	. 050	. 058	. 106	. 097	. 149	. 196	. 224	. 235
Solution Time	. 025	. 226	. 689	. 903	3.053	3.026	7.022	9.829	15, 180	14.622
OPHELIE/LP <sup>1</sup> Solution Time	.755	4.012		39, 375						276.90
Out-of-Kilter Solution Time	. 165	.617	1.248	1.748	3,312	3,639	6.152	11.903	14.587	12, 302
Density <sup>2</sup>	. 35	. 65	09.	. 36	. 54	. 20	. 28	. 31	. 28	. 20
Problem Size	10 × 10	20 × 20	30 x 30	40 x 40	50 x 50	09 × 09	70 x 70	80 x 80	06 × 06	100 × 100

<sup>1</sup>All times are median times with five problems per group.

<sup>2</sup>Median density.

\*Iteration time.

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Table 2

Total Solution Time Relative to Start Time For Nondense Problems Using the Primal Algorithm<sup>1</sup>

¥ K	8	8	900	8	8	980	8	8.	8	8	8	8	010	010	
lke Iteration Time	8	. 502	. 923	1.133	2. 035	2, 117	3, 783	5. 209	6.277	6. E24	11.953	14. 690	20, 492	23, 386	
No. of Her.		88	158	218	308	355	558	140	825	906	1425	1663	2154	2477	
Northwest Corner Rule ton Start No. of Ite	•10.	.031	110	.050	. 151	8	.092	. 113	141	. 146	. 220	. 252	346	. 312	
North Colution Time	. 057	. 590	1.033	1. 186	2, 186	2. 186	3, 875	5. 322	6.418	6.970	12. 173	14, 942	20, 838	23.698	
¥	8	80	8	80	8	8	8	88	8	8	.011	010	.011	.011	
nimum² Ite ration Time	\$	. 249	. 532	889	1. 239	1.417	2.009	3.100	3, 757	4. 228	6. 323	8. 539	11.947	13.472	
Row-Column Minimum <sup>2</sup> Start No. of Iteratio Time Iter. Time	•	8	82	110	195	228	249	378	480	505	28	200	1089	1205	
Row-Co Start Time	.07	. 033	. 051	. 067	. 077	.084	. 163	171	. 174	. 175	3	. 501	. 730	013.	
Solution	.061	. 282	. 583	. 755	1.316	1.501	2,252	3.271	3.931	4.403	6.871	9.040	12.677	14. 282	
ř.	.005	98	<b>500</b>	900	. 007		.8	80	.00	808	010	8	010.	.01	
m² Iteration Time	. 056	. 226	. 497	969	1, 189	1, 397	2.068	2.990	3, 521	4. 122	5. 558	7.907	11.041	13, 952	
Row Minimum <sup>2</sup> t No. of Iter e Iter. Ti	12	39	66	110	168	204	289	378	490	205	570	849	1059	1222	
Row Start Time	.018	.034	. 046	170.	960	. 092	. 129	176	. 176	180	. 299	.311	.377	8	
Solution	. 074	. 260	. 543	. 767	1.284	1.489	2, 197	3. 166	3.697	4. 302	5.857	8.218	11.418	14.443	
API	. 005	900	.00	900	8.	900	90	.87	900	800	010	.016	.011	010	
Vogel Approximation Method <sup>2</sup> ion Start No. of Iteration ite Time Iter. Time	. 025	. 117	. 356	. 393	. 838	. 748	1.219	1.999	2, 279	2.445	4. 322	5.627	6.270	9, 133	
rimation Na of Iter.	s,	22	\$	67	140	143	158	248	286	291	450	573	296	806	
el Appros Start Time	.045	. 137	. 258	.469	. 717	. 845	1.356	1.651	2.068	2.495	3.979	5.612	7, 165	9.038	
Solut	.070	. 254	. 624	. 862	1, 555	1.593	2, 575	3.650	4, 347	4.940	8.301	11, 239	13, 435	18, 171	
Density <sup>3</sup>	. 35	. 65	9	. 36	ă,	. 20	. 28	.31	. 28	. 20	. 37	.31	61 .	. 17	
Problem Size	10 × 10	20 × 20	30 x 30	40 × 40	50 x 50	09 × 09	70 × 70	80 x 80	06 × 06	130 × 100	125 x 125	150 x 150	175 x 175	200 x 200	•

The "modified first negative" basis charge criterion was used in testing the start algorithms.  $^2$ All times are median times with five problems per group.  $^3$ Median density.

Average per Iteration.

Table 3

Total Solution Times Relative to Start Times For Dense Problems Using the Primal Algorithm<sup>1</sup>

Northwest Corner Rule Solution Start No. of Bereston API* Time Time Iter. Time API*	.005 .143 .015 30 .128 .004	700 182, 68 500 618, 700	.006 1.253 .052 157 1.210 .008	.008 2.689 .079 282 2.610 009	.009 4.069 .110 366 3.959 .011	.010 6.441 .143 533 6.298 .012	.010 9.467 .170 693 9.297 .013	. 612 12.958 . 214 801 12.744 . 016	.013 17.533 .267 882 17.266 .017	.014 21.590 .320 1170 21.270 .018
Row-Column Minimum <sup>2</sup> tart Na of Heration 'Ime Iter, Time	14 .069	36 . 239	625 . 529	101 . 815	147 1.317	215 2.086	235 2. 423	327 3, 904	398 5, 123	410 5.635
Solution Start	109 040	. 328 . 089	. 670 . 141	1.027 . 212	1, 642 , 325	2, 491 , 405	2, 922 . 499	4.519 .615	5.845 .722	6.515 .880
 VH•	.003	<b>90</b> °.	.005	.00	.00	600	.010	.010	010.	110.
n <sup>2</sup>   Iteration   Time	. 055	. 173	.375	. 744	1.075	1.590	2.477	3, 326	3, 788	4.298
Row Minimum <sup>2</sup> art Na of me Iter.	91	#	82	100	151	181	240	348	379	386
r S	. 019	. 038	. 065	960	. 136	. 179	. 229	. 283	. 340	. 402
Solution Time	. 074	. 211	. 440	. 840	1,211	1, 769	2. 706	3.609	4.128	4. 700
on API	8	900	600	800	600	. o10	.010	.012	10.	. 015
Method Iteration Time	.061	. 131	. 232	.545	. 882	1, 353	1, 902	2,469	3.070	3, 312
oximation No of Iter.	=	21	27	9	66	133	186	209	227	223
Vogel Approximation Method Solution Start No. of Iteration Time Time Iter. Time	.051	. 165	. 339	. 556	. 885	1,217	1,647	2, 171	2.633	3, 257
Vol Solution Time	. 112	. 296	. 571	1. 101	1.767	2.570	3, 549	4.640	5. 703	6.569
Problem Size Dense	10 x 10	20 x 20	30 x 30	40 × 40	50 x 50	09 × 09	07 × 07	80 x 80	06 × 06	100 x 100

 $^1{\rm The}$  "modified first negative" basis criteria was used in testing the start algorithms.  $^2{\rm All}$  times are median times with ten problems per group.

Average per iteration time.

Total Solution Time Relative to Basis Change Criteria For Nondense Problems Using the Primal Algorithm

Table 4

	· idv	.00	. 011	.018	. 022	.029	820	648	3	. 082	8	31.	15.	190	
	t lter. Time	810	. 189	529	800	1, 794	27.2.2	4.083	4. 880	10, 386	10.965	32, 240	36, 780	54. 224	
	egative <sup>2</sup> No. of Iter. Iter. Time	•	11	62	36	ē	82	1	110	126	187	176	233	285	
	Most Negative Start No. of 11	25	138	. 295	166	. 750	808	1.211	1.544	2.063	2. 429	4.000	5.682	7, 295	
	Solution	. 063	. 327	. 824	1. 266	2, 544	3, 180	5. 354	6. 424	12, 469	13, 394	36, 240	42.462	61.519	
	A PI•	8	.005	8	80	900	800	8	.013	. 013	.013	. 025	016	.018	. 020
		. 027	. 120	408	465	884	806	1.473	2.082	2.612	3. 109	5. 736	7, 353	9.039	11.045
	Nega lo. of	-	22	51	20	112	115	156	184	198	242	226	452	510	555 1
	Modified Most Negative <sup>2</sup> Solution Start No. of Iter. Time Time Iter. Time	.034	. 135	.274	.471	. 651 1	. 883 1	1. 320 1	1.680 1	2.056 1	2.511.2	4.041 2	5, 283 4	7, 188 5	9.471 5
	Modifi Solution S Time T	18	. 255	. 632	936	1.505	1. 792	2. 793 1	3. 762 1	4.668 2	5. 620 2	9.777 4	12,636 5	16, 227 7	20.516 9
	ģĒ	٠.	•	•		1	-	7	က်	÷	'n	Ġ	12.	16.	20.
	API*	. 005	900	88.	900	88	8	.007	900	800	800	010	010	.011	010
1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1		. 025	. 117	. 356	. 393	. 838	. 748	1.219	1.999	2.279	2,445	4.322	5.627	6.270	9, 133 .010
	rst Negative No. of Iter. Iter. Tim	•	23	9	67	140	143	158	248	286	291	450	573	296	803
		. 945	. 137	. 258	. 469	7117.	. 845	1, 356	1.651	2.068	2.495	3.979	5.612	7. 165	9.038
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	API.	900	900	80.	\$00.	900	900	900	900	700	100	600	010	600	606
	2 Iter. Time API*	. 028 . 004	. 125 . 005	. 367 . 005	.458 ,005	. 704 . 005	1.142 .005	1.317 .006	2, 124 , 006	2.908 .007	3.088 .007	4.890 .009	6.661 .010	8.653 .009	11.875
	Negative <sup>2</sup> No of Iter. Iter. Time	-	25	73	6	137	219	231	334	00+	446	261	693		1265 1
	First Negative Start No. of It	. 038	. 149	. 219	.455	722	. 920	1.261	1.677	2.061	2, 382	4.012	5.642	7, 205 1002	9,178 1265 11.875 .009
	Solution	990	. 274	. 586	. 913	1, 426	2.062	2,578	3.801	4.969	5.470	8.902	12, 303	15.858	21.053
	Density	. 35	. 65	09	. 36	. 54	20	. 28	. 31	. 28	. 20	. 37	.31	. 13	. 17
	Problem Size D	10 × 10	20 x 26	30 x 30	40 x 40	50 x 50	09 × 09	70 × 70	80 × 80	06 × 06	100 × 100 . 20	125 x 125	150 x 150 . 31	175 × 175 . 19	200 x 200 . 17

<sup>1</sup>The VAM start method was used for determinus an initial primal feasible basic solution.

<sup>2</sup>All times are median times with five problems per group.

<sup>3</sup>Median density.

\*Average per Iteration.

Table 5

Total Solution Time Relative to Basis Change Criteria For Dense Problems Using the Primal Algorithm<sup>1</sup>

*FI	800	. 017	. 034	950	. 063	. 118	. 157	. 203	.257	
Iter. Time	3	. 292	.673	2, 260	3.896	8.580	14.271	24.767	30, 916	
Most Negative <sup>2</sup> Start No. of Fime Iter.	•	11	20	7	4.	(T)	16	122	121	
Most N. Start Time	\$	. 162	. 331	. 561	. 883	1.30	1.648	2, 160	2.640	
Solution	. 133	. 488	1.022	2.851	4.861	908.6	15.978	26.988	33.614	
API.	3	8	.011	. 016	.016	920	. 022	.027	. 032	. 036
re <sup>2</sup> Reration Time	. 050	191	. 342	. 805	1.201	2. 138	2. 524	4.069	6.032	5.350
Negativ No of Iter.	10	23	8	57	5	103	122	149	222	9
Modified Most Negative <sup>2</sup> ution Start No. of 1te me Time Iter.	. 948	. 154	. 320	.573	2	1, 199	1.682	2.170	2.642	3, 343
Modil Solution Time	. 111	. 359	889	1, 384	2.072	3,365	4.239	6.281	3.806	9.119
AM.	8	8	600	808	600	010	. 010	. 012	<b>10</b> .	. 015
teration Time	.081	. 131	. 232	. 545	. 882	1, 353	1.902	2,469	3.070	3, 312
Negativ No of Iter.	<b>±</b>	21	27	9	66	133	186	<b>308</b>	227	223
Modified First Negative <sup>2</sup> lution Start No. of It 'lme Time Iter.	.051	. 165	. 339	. 556	. 885	1.217	1,647	2.171	2.633	3,257
Modif Solution Time	. 112	. 296	.571	1.101	1.767	2.570	3.549	4.640	5. 703	6.569
<b>AM</b> •	8	80.	800	.007	800	8	010	60	. 013	.012
ter. Time	.042	. 225	. 375	2	1.055	1.595	2.256	2.977	4.171	4. 129
First Negative <sup>2</sup> Start No. of Time Iter.	01	37	9	90	125	173	227	318	331	329
First N Start Time	. 050	. 161	. 333	. 558	. 864	1.212	1,654	2, 159	2.646	3, 335
Solution Time	. 092	. 386	. 708	1.202	1.919	2.807	3.910	5, 136	6.817	7.464
Problem Size Dense	10 × 10	20 x 20	30 x 30	40 x 40	50 x 50	90 × 90	70 x 70	80 x 80	06 × 06	100 x 100

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 $^1{\rm The~VAM}$  start method was used to determine an initial primal feasible basic solution.  $^2{\rm All}$  times are median times with ten problems per group.

\*Average per Iteration.

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indicated in Table 1, the out-of-kilter algorithm was more efficient than the special dual a gorithm (but not the primal algorithm) as soon as m became at least 80.

To compare the change of basis criteria the same starting solution was used. Consequently, only the main loop times required comparison to isolate the most effective criterion. The data in Tables 4 and 5 indicate that the highly popular and often referenced "most negative" criterion for determining the next vector to enter the basis is not the most efficient basis change criterion. In fact, as indicated in Tables 4 and 5, for the basis change criteria examined in the experiment, the most negative evaluator criterion was consistently slower than the other basis change criteria examined. Furthermore, when m became larger than 100, the "most negative" criterion required at least three times as much computational effort as the other basis change criteria.

Interestingly, the data in Tables 1 and 4 indicate that the out-of-kilter algorithm is more efficient computationally than the primal row-column sum algorithm, when the latter employs the most negative evaluator criterion. Perhaps the folklore concerning the superiority of the out-of-kilter method was developed by comparing the out-of-kilter network algorithm to a transportation method that used the most negative basis change criterion. (The lack of techniques such as those indicated in Section 2.1 in the previously developed special purpose primal methods would undoubtedly have accentuated the apparent difference between primal and out-of-kilter methods.)

The data in Tables 2 and 3 (using the row-minimum start) support the finding of Dennis [10] that the total solution time required to solve m x m distribution problems increases proportionately to m<sup>2</sup>. Also, given the same relative cost range for a distribution problem, the data indicate that the total solution time is approximately twice the time necessary to obtain an initial feasible basic solution for the VAM start in dense and nondense problems (see Tables 2 and 3).

# 3.2 Solution Algorithms, Start Methods and Basis Change Criteria

An important factor influencing total computation time (and hence computational efficiency) is the interrelationship between the start methods and basis change criteria. The relevant tradeoffs for the start algorithms involve the time required for obtaining an initial feasible basic solution and the number of iterations subsequently required to complete the solution of the problem. Similarly, the relevant tradeoffs for the basis change criteria involve the time consumed in searching for a new vector to enter the basis and the number of iterations required to obtain an optimal solution (time per iteration versus total number of iterations).

To determine the effect of combining various start procedures with various change of basis criteria, and to avoid excessive computational expense in making this determination, we tested each change of basis criterion relative to a single start method (VAM), and then tested each of the start methods relative to the change of basis criterion that was found to be best.

Ideally, had it not involved an enormous expenditure of computer time, it would have been desirable to have tested all combinations of start procedures and change of basis criteria against each other. However, our findings were sufficiently clear cut to make it a reasonably safe bet that the change of basis criterion and start procedure that were found to be best "in isolation" will also be best in combination. In particular, from Table 4 and Table 5 it is clear that the "modified first negative" basis is faster than the other criteria examined both in terms of the time per iteration (or "main loop time") and in terms of total solution time. (This result concurs with the finding of Dennis [10].)

In general, the number of iterations required to obtain the optimal solution was significantly reduced by using the "modified most negative" and "most negative" basis change criteria. As indicated in Tables 4 and 5, these basis change criteria led to a total iteration count that was consistently less than 80 percent of the iteration count associated with the "first negative" and "modified first negative" basis change criteria.

More importantly, however, the "first negative" and "modified first negative" basis change criteria are faster than the other criteria examined both in terms of the time per iteration (or "main loop time") and in terms of total time. The efficiency of the "first negative" and "modified first negative" criteria is accentuated as m increases and the density of the distribution problem approaches 100%. The data in Tables 4 and 5 indicate that the main loop time for the "first negative" and the "modified first negative"

basis change criteria is consistently less than 70% of the main loop time required by the "modified most negative" and "most negative" basis change criteria. Furthermore, these data indicate that the "modified first negative evaluator" criterion consistently requires 20 percent fewer iterations than the "first negative evaluator" criterion and the main loop time of the two criteria were approximately equal. Consequently, as indicated by the data in Tables 4 and 5, the "modified first negative" criterion was consistently at least 15 percent faster in terms of total computation time than the "first negative" criterion.

Another interesting fact that the study revealed about the basis change criteria concerns the number of possible nonbasic variables that are examined during each iteration. The "first negative" and "modified first negative" criteria always evaluate substantially fewer nonbasic variables than the "most negative" criterion, and as m increases the difference in the number of nonbasic variables evaluated also increases significantly. The data in Tables 4 and 5 indicate that when m equals 20 and 200, the number of nonbasic variables that must be examined by the "first negative" and "modified first negative" criteria as compared to the "most negative" criterion is only 50% and 5%, respectively. Thus the results indicate that the "modified first negative" basis change criterion provides the best compromise of the tradeoffs for both dense and nondense problems.

Having determined the "modified first negative" change of basis criterion to be significantly more effective than the others, we used this

choice rule uniformly thereafter in conducting the tests of the alternative start methods. Again our results were clear cut. From Tables 2 and 3 it is apparent that the Row Minimum start method is best for both dense and nondense problems.

There are a number of interesting observations to be obtained from Tables 2 and 3. Vogel's method, while yielding a good start from the standpoint of the number of iterations required after it is completed, takes an inordinate amount of time to find an initial solution, and exhibits a large average iteration time for dense problems, indicating that the loop structure is complicated. (See Table 2.) Similarly, the loop structure for the Northwest Corner start is complicated, whereas the loop structures for the Row Minimum and Row-Column Minimum are simpler. On dense problems the solution time of VAM and Row-Column Minimum are approximately equal.

Comparing the results in Tables 2 and 3 yields several insights into the effect of density on primal codes. In general the nondense problems were solved only about 10% faster than dense problems using the Row Minimum start, and, in fact, the dense problems were solved faster for problem sizes not exceeding 50 x 50. Thus density does not appreciably affect total solution time. Average main loop time is substantially reduced by density, indicating that loop structure becomes less complicated as the density of a problem decreases. This reduction in average main loop time is partially offset by the fact that the number of iterations increases slightly as density decreases.

As indicated in Tables 1 and 2, the special purpose primal simplex algorithm was consistently faster than the special purpose dual simplex algorithm. It would be premature, however, to conclude that one should abandon the development of special purpose dual simplex algorithms for a number of reasons. For instance, this study indicates that unlike the primal approach, the dual approach is greatly affected by density. A 100 x 100 problem of 13% density was solved in 7.697 seconds while a 100 x 100 problem of 20% density was solved in 14.622 seconds. Although 7.697 seconds is still nearly twice the median solution time for 100 x 100 problems using the primal approach, it does suggest that the dual method may be an efficient solution procedure for solving sparse network problems. It is interesting to note that the out-of-kilter code was not nearly as sensitive to density, since its times on the above problems were 9.973 and 12.045 seconds, respectively, hence making the dual method appear still more attractive for sparse problems. Further support for this possibility is given by noting that the solution time of the primal approach varies proportionately to the number of nodes in the problem while the dual approach varies proportionately to the number of arcs in the problem.

The dual start method used in this study was not an advanced start procedure. Even so, the special purpose dual simplex algorithm consistently required fewer than 50% of the number of iterations necessary for the primal simplex algorithm when any start method except VAM was used. These observations indicate that substantial computational saving may be present if

advanced start algorithms (such as those developed in [16]) are employed in conjunction with the special purpose dual simplex algorithm.

Although the dual algorithm appears to be computationally slower than the primal algorithm, it is particularly useful as compared to the primal algorithm for conducting certain types of postoptimality analyses. A principal motivation cited in the literature for applying the dual method to distribution problems (see, e.g., [2,3,7]) is the possibility that the problem's "supplies" and "demands" may not be permanently fixed, but rather are subject to change. In such a situation, the ability to begin from an optimal basis to a given problem and proceed via the dual method to an optimal solution for the altered problem is extremely useful. Such an approach is not available using the primal algorithm without extensive computation.

# 4.0 Concluding Remarks

In the context of the important class of linear programs known as distribution (or transportation) problems, the special purpose primal and dual simplex and the out-of-kilter algorithms have been shown to be consistently faster than the general simplex algorithm. The special purpose primal simplex algorithm is in turn consistently faster than the out-of-kilter algorithm and the dual simplex algorithm. Our results indicate that it is important to determine a "good" starting basis for the dual algorithm in order to obtain an advantageous tradeoff between the total number of iterations and the time required to implement the starting algorithm.

The computational efficiency of the special purpose primal and dual algorithms for solving distribution problems has been shown to be particularly sensitive to the search time required to find a new arc to enter the basis at each iteration. The dual algorithm is more sensitive in this regard than the primal algorithm. The relative slowness of the dual algorithm's main loop time suggests that further research on advanced dual "start" methods is worthwhile, since a relatively large amount of time could be spent finding an initial basic feasible solution and still permit an overall reduction in computation time by reducing the number of iterations.

On the other hand, a highly efficient dual start will not succeed in making the special purpose dual algorithm faster than the special purpose primal algorithm unless the main loop time for the dual method can be greatly reduced. An important qualification to this assertion rests on the fact that the special purpose dual algorithm appears to gain efficiency relative to the primal algorithm as the proportion of the number of nodes to number of admissible cells increases. This, combined with the fact that a basic feasible dual solution exists for networks (see [19]), indicates that the special purpose dual algorithm might be the best solution procedure for minimum cost network problems. <sup>5</sup>

In the domain of choice rules, the data indicate that the number of problem-solving iterations is substantially reduced by using the "most negative" and "modified most negative basis change criteria. However, this reduced iteration count does not produce a correspondingly reduced computation time since the search time required by these criteria (particularly for dense problems) is substantially greater for the "first negative" and "modified first negative"

criteria. In consequence, the latter criteria turn out to be more attractive from the standpoint of total computation time. The "modified first negative" criterion is generally the more efficient of the two, although for non-dense problems the "first negative" and "modified first negative" criteria appear to be comparably efficient.

The most efficient form of the special primal simplex algorithm was found to be based on the Row Minimum start algorithm and the "modified first negative" basis change criterion. The primal simplex algorithm employed in this study proved to be from 2 to 15 times faster than the other special purpose distribution algorithms that were available for examination.

In evaluating the performance of the methods tested, considerable effort was made to insure that the computer codes were comparably efficient in implementing the underlying algorithms, start methods, and basis change criteria. Moreover, the large differences found between the performance of the various procedures tested leads us to believe that these differences will continue to hold for improved computer codes (e.g., codes that take special advantage of the characteristics of the computer on which they are implemented). In fact the techniques used to time the codes suggest that the relative superiority of the special purpose algorithms may be even greater than indicated in Section 3. Specifically, the codes for the special purpose methods included a number of clocks to measure different aspects of the solution process, whereas the codes for the general simplex algorithms and the out-of-kilter algorithm only included clocks to measure the total solution time. When the

clocks are removed from the special purpose computer codes, the computational efficiency of the algorithms has been found to increase by as much as 40 percent. 6

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## **FOOTNOTES**

 $^1$ It is interesting to note that most of the folklore related to the computational requirements for solving distribution problems is based on computation performed 10 to 15 years ago. Furthermore, the distribution problems examined were often small, maximum dimension of 20 x 20 [6,9], and they were solved either by hand or by first or second generation computers.

- <sup>2</sup> This code was developed by F. Glover, D. Karney and D. Klingman.
- $^{3}$ This code was developed by the authors.
- 4"Main loop time" (average iteration time) is the time required to test for optimality to determine the variable to enter the basis, to obtain the variable to leave the basis, to find the basis exchange path, and to update the basic variables.

<sup>5</sup>The authors are currently testing other dual start algorithms and new procedures for reducing the main loop time of the dual method. Network algorithms are also being tested.

<sup>6</sup>Incidental to our investigation was the finding--or rather confirmation--that one can make no valid conjecture about the computational efficiency of two implementations of the same algorithm unless the two computer codes are executed on the same machine using the same data and the same compiler. This fact was underscored in tests comparing our special purpose primal simplex computer code to another special purpose code, based on similar techniques, that was executed on a Univac 1108. The latter code ran faster on the Univac than our code ran on the CDC 6600, suggesting that our code was less efficient due to the fact that the CDC 6600 is generally a faster machine than the 1108.

Surprisingly, when the Univac 1108 computer code was executed on the CDC 6600, it ran slower than the special purpose primal computer code used in this experiment. The apparent reason for this reversal is that the Univac 1108 FORTRAN compiler produces more efficient object code.

This result stimulated us to check with Control Data Corporation about their compiler, which led us to discover that the CDC 6600 has two FORTRAN compilers. One produces an object code which executes about twice as fast as the other. This implies that our times could have been reduced by 50% if we had used the other compiler. Since the total compilation and execution time of the two compilers is equal, The University of Texas Computation Center only uses the less efficient object code compiler.

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